

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061925\
 Data File : BN037337.D
 Acq On : 19 Jun 2025 05:54
 Operator : RC/JU
 Sample : PB168508BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB168508BS

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 06/19/2025
 Supervised By :mohammad ahmed 06/23/2025

Quant Time: Jun 19 06:28:28 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN061925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 18 18:40:16 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.568	152	1288	0.400	ng	0.00
7) Naphthalene-d8	10.340	136	2690	0.400	ng	#-0.01
13) Acenaphthene-d10	14.213	164	1686	0.400	ng	-0.01
19) Phenanthrene-d10	16.971	188	3212	0.400	ng	0.00
29) Chrysene-d12	21.171	240	3029	0.400	ng	# 0.00
35) Perylene-d12	23.354	264	3232	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.170	112	825	0.352	ng	0.00
5) Phenol-d6	6.752	99	802	0.348	ng	0.00
8) Nitrobenzene-d5	8.717	82	760	0.339	ng	-0.01
11) 2-Methylnaphthalene-d10	11.945	152	1622m	0.371	ng	0.00
14) 2,4,6-Tribromophenol	15.718	330	329	0.265	ng	0.00
15) 2-Fluorobiphenyl	12.838	172	2716	0.368	ng	0.00
27) Fluoranthene-d10	19.008	212	3351	0.329	ng	0.00
31) Terphenyl-d14	19.616	244	2523	0.363	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.104	88	501	0.436	ng	# 44
3) n-Nitrosodimethylamine	3.415	42	455	0.341	ng	90
6) bis(2-Chloroethyl)ether	7.004	93	760	0.386	ng	100
9) Naphthalene	10.394	128	2428	0.349	ng	# 92
10) Hexachlorobutadiene	10.682	225	1119	0.318	ng	# 99
12) 2-Methylnaphthalene	12.021	142	1521	0.310	ng	96
16) Acenaphthylene	13.935	152	2681	0.378	ng	98
17) Acenaphthene	14.277	154	1575	0.334	ng	99
18) Fluorene	15.271	166	2259	0.338	ng	98
20) 4,6-Dinitro-2-methylph...	15.378	198	302	0.306	ng	89
21) 4-Bromophenyl-phenylether	16.177	248	914	0.343	ng	96
22) Hexachlorobenzene	16.276	284	1066	0.382	ng	93
23) Atrazine	16.450	200	743	0.357	ng	# 89
24) Pentachlorophenol	16.624	266	250	0.162	ng	94
25) Phenanthrene	17.009	178	3326	0.368	ng	99
26) Anthracene	17.108	178	3130	0.369	ng	100
28) Fluoranthene	19.040	202	4162	0.331	ng	99
30) Pyrene	19.402	202	4189	0.363	ng	99
32) Benzo(a)anthracene	21.153	228	3675	0.377	ng	99
33) Chrysene	21.206	228	4517	0.363	ng	99
34) Bis(2-ethylhexyl)phtha...	21.099	149	1537	0.384	ng	98
36) Indeno(1,2,3-cd)pyrene	25.541	276	5500	0.405	ng	# 93
37) Benzo(b)fluoranthene	22.702	252	4039	0.358	ng	98
38) Benzo(k)fluoranthene	22.746	252	4639	0.374	ng	96
39) Benzo(a)pyrene	23.260	252	4040	0.386	ng	96
40) Dibenzo(a,h)anthracene	25.555	278	4089	0.375	ng	# 88
41) Benzo(g,h,i)perylene	26.204	276	4647	0.373	ng	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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