

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100923\
 Data File : BN028160.D
 Acq On : 09 Oct 2023 14:34
 Operator : MA/JU
 Sample : PB155953BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB155953BS

Quant Time: Oct 09 15:29:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 09 12:40:55 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	3591	0.400	ng	0.00
7) Naphthalene-d8	10.746	136	10149	0.400	ng	# 0.01
13) Acenaphthene-d10	14.566	164	4710	0.400	ng	0.00
19) Phenanthrene-d10	17.331	188	8895	0.400	ng	0.00
29) Chrysene-d12	21.515	240	8044	0.400	ng	0.00
35) Perylene-d12	23.984	264	8695	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.473	112	3576	0.334	ng	0.00
5) Phenol-d6	7.091	99	4152	0.307	ng	0.00
8) Nitrobenzene-d5	9.123	82	3486	0.404	ng	0.00
11) 2-Methylnaphthalene-d10	12.328	152	7050	0.469	ng	0.00
14) 2,4,6-Tribromophenol	16.077	330	524	0.243	ng	0.00
15) 2-Fluorobiphenyl	13.197	172	7874	0.464	ng	0.00
27) Fluoranthene-d10	19.360	212	8971	0.402	ng	0.00
31) Terphenyl-d14	19.946	244	8115	0.432	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.357	88	1547	0.384	ng	# 76
3) n-Nitrosodimethylamine	3.704	42	2010	0.436	ng	# 94
6) bis(2-Chloroethyl)ether	7.366	93	3877	0.356	ng	92
9) Naphthalene	10.789	128	11044	0.421	ng	99
10) Hexachlorobutadiene	11.013	225	2009	0.431	ng	# 99
12) 2-Methylnaphthalene	12.404	142	6437	0.351	ng	98
16) Acenaphthylene	14.298	152	8059	0.386	ng	99
17) Acenaphthene	14.630	154	5459	0.410	ng	99
18) Fluorene	15.618	166	6633	0.384	ng	98
20) 4,6-Dinitro-2-methylph...	16.785	198	45	0.381	ng	84
21) 4-Bromophenyl-phenylether	16.512	248	1787	0.385	ng	# 83
22) Hexachlorobenzene	16.599	284	2142	0.440	ng	94
23) Atrazine	16.785	200	1349	0.333	ng	95
24) Pentachlorophenol	16.971	266	6769	2.954	ng	94
25) Phenanthrene	17.368	178	9882	0.412	ng	99
26) Anthracene	17.468	178	8252	0.364	ng	99
28) Fluoranthene	19.388	202	10910	0.394	ng	99
30) Pyrene	19.755	202	11393	0.386	ng	98
32) Benzo(a)anthracene	21.506	228	10609	0.388	ng	99
33) Chrysene	21.560	228	11333	0.430	ng	99
34) Bis(2-ethylhexyl)phtha...	21.372	149	4605	0.256	ng	100
36) Indeno(1,2,3-cd)pyrene	26.568	276	14610	0.500	ng	99
37) Benzo(b)fluoranthene	23.218	252	11920	0.424	ng	98
38) Benzo(k)fluoranthene	23.271	252	11368	0.399	ng	96
39) Benzo(a)pyrene	23.873	252	10209	0.384	ng	95
40) Dibenzo(a,h)anthracene	26.589	278	11855	0.496	ng	97
41) Benzo(g,h,i)perylene	27.378	276	12660	0.516	ng	98

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

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