

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
 Data File : BN033155.D
 Acq On : 01 Aug 2024 06:31
 Operator : MA/JU
 Sample : P3265-12MSD
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4D42MSD

Quant Time: Aug 01 07:41:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 14:49:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.520	152	2199	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.273	136	7580	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.138	164	4380	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.911	188	9722	0.400	ng/ul	-0.03
17) Chrysene-d12	21.128	240	8811	0.400	ng/ul	-0.02
23) Perylene-d12	23.308	264	9197	0.400	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.056	96	1355	0.487	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.868	152	3621	0.293	ng/ul	-0.06
18) Fluoranthene-d10	18.944	212	8706	0.349	ng/ul	-0.03
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.086	88	3778	1.191	ng/ul#	71
5) Naphthalene	10.322	128	6501	0.334	ng/ul	98
7) 2-Methylnaphthalene	11.945	142	4216	0.320	ng/ul	99
8) 1-Methylnaphthalene	12.165	142	4253	0.298	ng/ul	99
10) Acenaphthylene	13.870	152	6575	0.367	ng/ul	98
11) Acenaphthene	14.203	153	5401	0.401	ng/ul	99
12) Fluorene	15.216	166	5783	0.368	ng/ul	99
14) Pentachlorophenol	16.599	266	1045	0.528	ng/ul	97
15) Phenanthrene	16.954	178	24442	0.937	ng/ul	99
16) Anthracene	17.051	178	10336	0.441	ng/ul	99
19) Fluoranthene	18.977	202	48713	1.679	ng/ul	98
20) Pyrene	19.348	202	46418	1.492	ng/ul	95
21) Benzo(a)anthracene	21.110	228	27355	0.910	ng/ul	98
22) Chrysene	21.166	228	28568	0.815	ng/ul	99
24) Benzo(b)fluoranthene	22.656	252	34437	1.282	ng/ul	90
25) Benzo(k)fluoranthene	22.697	252	18843	0.585	ng/ul#	90
26) Benzo(a)pyrene	23.217	252	23603	0.933	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	25.500	276	22231	0.578	ng/ul	97
28) Dibenzo(a,h)anthracene	25.493	278	10267	0.337	ng/ul	98
29) Benzo(g,h,i)perylene	26.178	276	19982	0.654	ng/ul	94

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

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