

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121824\
 Data File : BN035695.D
 Acq On : 19 Dec 2024 06:42
 Operator : RC/JU
 Sample : PB165622BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS622

Quant Time: Dec 19 08:33:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN121624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 16:26:08 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.276	152	2854	0.400	ng/ul	0.00
4) Naphthalene-d8	10.020	136	7609	0.400	ng/ul	-0.01
9) Acenaphthene-d10	13.940	164	5215	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.713	188	9797	0.400	ng/ul	0.00
17) Chrysene-d12	20.955	240	8993	0.400	ng/ul	0.00
23) Perylene-d12	23.039	264	9867	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	1614	0.638	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.626	152	4022	0.371	ng/ul	-0.01
18) Fluoranthene-d10	18.763	212	8740	0.333	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.984	88	4417	1.729	ng/ul#	71
5) Naphthalene	10.069	128	7030	0.377	ng/ul	99
7) 2-Methylnaphthalene	11.703	142	4306	0.343	ng/ul	99
8) 1-Methylnaphthalene	11.928	142	4545	0.349	ng/ul	99
10) Acenaphthylene	13.653	152	6649	0.329	ng/ul	98
11) Acenaphthene	14.000	153	5143	0.337	ng/ul	100
12) Fluorene	15.008	166	5503	0.310	ng/ul	98
14) Pentachlorophenol	16.375	266	1481	0.590	ng/ul	98
15) Phenanthrene	16.751	178	8369	0.341	ng/ul	100
16) Anthracene	16.844	178	7311	0.319	ng/ul	100
19) Fluoranthene	18.795	202	10124	0.315	ng/ul	100
20) Pyrene	19.163	202	10725	0.303	ng/ul	99
21) Benzo(a)anthracene	20.941	228	9967	0.334	ng/ul	100
22) Chrysene	20.991	228	11124	0.359	ng/ul	99
24) Benzo(b)fluoranthene	22.428	252	10251	0.353	ng/ul	97
25) Benzo(k)fluoranthene	22.472	252	11173	0.347	ng/ul	97
26) Benzo(a)pyrene	22.951	252	9863	0.350	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.067	276	14107	0.411	ng/ul#	52
28) Dibenzo(a,h)anthracene	25.084	278	11351	0.434	ng/ul#	94
29) Benzo(g,h,i)perylene	25.681	276	10203	0.371	ng/ul	96

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

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