

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
 Data File : BN033128.D
 Acq On : 31 Jul 2024 12:48
 Operator : MA/JU
 Sample : P3265-03DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DQ8DL

Quant Time: Jul 31 13:21:45 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	7.528	152	1993	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.300	136	7068	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.152	164	4655	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.928	188	10592	0.400	ng/ul	-0.01
17) Chrysene-d12	21.148	240	10945	0.400	ng/ul	0.00
23) Perylene-d12	23.334	264	11338	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.060	96	142	0.056	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.950	152	307	0.027	ng/ul	0.02
18) Fluoranthene-d10	18.972	212	1051	0.034	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.350	128	15979	0.880	ng/ul	96
10) Acenaphthylene	13.889	152	5771	0.303	ng/ul	97
11) Acenaphthene	14.217	153	13392	0.935	ng/ul	99
12) Fluorene	15.240	166	6385	0.382	ng/ul	99
15) Phenanthrene	16.971	178	7756	0.273	ng/ul	99
16) Anthracene	17.068	178	35468	1.389	ng/ul	97
19) Fluoranthene	19.000	202	15182	0.421	ng/ul	99
20) Pyrene	19.367	202	7510	0.194	ng/ul#	93
21) Benzo(a)anthracene	21.137	228	14002	0.375	ng/ul	99
22) Chrysene	21.183	228	7921	0.182	ng/ul	99
24) Benzo(b)fluoranthene	22.682	252	12188	0.368	ng/ul	99
25) Benzo(k)fluoranthene	22.723	252	19074	0.481	ng/ul	95
26) Benzo(a)pyrene	23.261	252	4596	0.147	ng/ul#	67
27) Indeno(1,2,3-cd)pyrene	25.544	276	23822	0.503	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.534	278	17280	0.460	ng/ul	94
29) Benzo(g,h,i)perylene	26.221	276	14564	0.387	ng/ul	96

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

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