

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
 Data File : BN033129.D
 Acq On : 31 Jul 2024 13:24
 Operator : MA/JU
 Sample : P3283-03DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DQ9DL

Quant Time: Jul 31 13:53:43 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	7.537	152	1768	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.306	136	6278	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.148	164	4298	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.920	188	9695	0.400	ng/ul	-0.02
17) Chrysene-d12	21.151	240	9864	0.400	ng/ul	0.00
23) Perylene-d12	23.337	264	10619	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.056	96	122	0.055	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.945	152	321	0.031	ng/ul	0.02
18) Fluoranthene-d10	18.963	212	1149	0.041	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.355	128	17362	1.077	ng/ul	96
10) Acenaphthylene	13.879	152	20574	1.170	ng/ul	98
11) Acenaphthene	14.213	153	31384	2.372	ng/ul	99
12) Fluorene	15.230	166	17475	1.133	ng/ul	99
15) Phenanthrene	16.962	178	49929	1.920	ng/ul	98
16) Anthracene	17.063	178	42417	1.814	ng/ul	97
19) Fluoranthene	18.995	202	40243	1.239	ng/ul	97
20) Pyrene	19.362	202	18537	0.532	ng/ul	100
21) Benzo(a)anthracene	21.140	228	4668	0.139	ng/ul	98
22) Chrysene	21.186	228	18457	0.470	ng/ul	99
24) Benzo(b)fluoranthene	22.682	252	15717	0.507	ng/ul	94
25) Benzo(k)fluoranthene	22.726	252	16225	0.436	ng/ul	95
26) Benzo(a)pyrene	23.247	252	21904	0.750	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	25.540	276	23496	0.529	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.530	278	22180	0.630	ng/ul#	91
29) Benzo(g,h,i)perylene	26.218	276	16915	0.480	ng/ul	96

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

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