

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
 Data File : BN033161.D
 Acq On : 01 Aug 2024 10:08
 Operator : MA/JU
 Sample : PB162364BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS364

Quant Time: Aug 01 11:01:12 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	7.507	152	2779	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.273	136	9972	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.143	164	5908	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.920	188	12567	0.400	ng/ul	-0.02
17) Chrysene-d12	21.140	240	10656	0.400	ng/ul	#-0.01
23) Perylene-d12	23.323	264	11487	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.052	96	2403	0.684	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.884	152	4993	0.307	ng/ul	-0.04
18) Fluoranthene-d10	18.963	212	10959	0.363	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.081	88	7229	1.804	ng/ul#	67
5) Naphthalene	10.322	128	9121	0.356	ng/ul	97
7) 2-Methylnaphthalene	11.961	142	6047	0.349	ng/ul	100
8) 1-Methylnaphthalene	12.165	142	6092	0.325	ng/ul	99
10) Acenaphthylene	13.870	152	8464	0.350	ng/ul	99
11) Acenaphthene	14.208	153	6926	0.381	ng/ul	100
12) Fluorene	15.221	166	7230	0.341	ng/ul	99
14) Pentachlorophenol	16.616	266	1252	0.489	ng/ul	98
15) Phenanthrene	16.962	178	11012	0.327	ng/ul	99
16) Anthracene	17.068	178	9721	0.321	ng/ul	100
19) Fluoranthene	18.995	202	13743	0.392	ng/ul	99
20) Pyrene	19.362	202	14444	0.384	ng/ul	98
21) Benzo(a)anthracene	21.125	228	11741	0.323	ng/ul	99
22) Chrysene	21.177	228	14658	0.346	ng/ul	99
24) Benzo(b)fluoranthene	22.671	252	12192	0.363	ng/ul	92
25) Benzo(k)fluoranthene	22.715	252	14442	0.359	ng/ul#	91
26) Benzo(a)pyrene	23.238	252	11863	0.376	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.530	276	15103	0.314	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.517	278	11798	0.310	ng/ul	99
29) Benzo(g,h,i)perylene	26.198	276	11946	0.313	ng/ul	96

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

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