

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080124\
 Data File : BN033165.D
 Acq On : 01 Aug 2024 12:32
 Operator : MA/JU
 Sample : PB162172BS
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS172

Quant Time: Aug 01 13:17:42 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	2175	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.295	136	7613	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.152	164	4549	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.933	188	9215	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	6945	0.400	ng/ul	0.00
23) Perylene-d12	23.337	264	7376	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.056	96	2014	0.732	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	3667	0.295	ng/ul	-0.02
18) Fluoranthene-d10	18.972	212	8067	0.410	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.086	88	5730	1.827	ng/ul#	68
5) Naphthalene	10.350	128	6750	0.345	ng/ul	99
7) 2-Methylnaphthalene	11.983	142	3879	0.293	ng/ul	96
8) 1-Methylnaphthalene	12.181	142	4132	0.289	ng/ul	100
10) Acenaphthylene	13.884	152	6043	0.325	ng/ul	98
11) Acenaphthene	14.217	153	5101	0.364	ng/ul	100
12) Fluorene	15.240	166	5215	0.320	ng/ul	98
14) Pentachlorophenol	16.637	266	713	0.380	ng/ul	100
15) Phenanthrene	16.979	178	7776	0.315	ng/ul	98
16) Anthracene	17.085	178	6464	0.291	ng/ul	98
19) Fluoranthene	19.005	202	9565	0.418	ng/ul	99
20) Pyrene	19.372	202	9967	0.406	ng/ul	98
21) Benzo(a)anthracene	21.139	228	7314	0.309	ng/ul	99
22) Chrysene	21.186	228	9845	0.356	ng/ul	100
24) Benzo(b)fluoranthene	22.685	252	7261	0.337	ng/ul	96
25) Benzo(k)fluoranthene	22.726	252	9685	0.375	ng/ul	96
26) Benzo(a)pyrene	23.252	252	7543	0.372	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.547	276	10922	0.354	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.540	278	8516	0.348	ng/ul	94
29) Benzo(g,h,i)perylene	26.225	276	7669	0.313	ng/ul	98

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

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