

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015430.D
 Acq On : 09 Jul 2021 04:52
 Operator : CG/JU
 Sample : PB137438BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS438

Quant Time: Jul 09 05:33:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.715	152	2927	0.400	ng/ul	0.00
4) Naphthalene-d8	10.479	136	9844	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.339	164	5055	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	10272	0.400	ng/ul	0.00
17) Chrysene-d12	21.284	240	8196	0.400	ng/ul	0.00
23) Perylene-d12	23.523	264	7097	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	2463	0.747	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	5458	0.353	ng/ul	0.00
18) Fluoranthene-d10	19.117	212	11223	0.417	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.315	88	6455	1.811	ng/ul	89
5) Naphthalene	10.529	128	10873	0.346	ng/ul	99
7) 2-Methylnaphthalene	12.146	142	7065	0.344	ng/ul	99
8) 1-Methylnaphthalene	12.366	142	6614	0.333	ng/ul	99
10) Acenaphthylene	14.057	152	8492	0.325	ng/ul	99
11) Acenaphthene	14.400	153	7021	0.351	ng/ul	99
12) Fluorene	15.390	166	7815	0.353	ng/ul	99
14) Pentachlorophenol	16.740	266	1962	0.863	ng/ul	99
15) Phenanthrene	17.128	178	13672	0.367	ng/ul	100
16) Anthracene	17.217	178	11335	0.345	ng/ul	99
19) Fluoranthene	19.150	202	14842	0.392	ng/ul	97
20) Pyrene	19.512	202	15021	0.395	ng/ul	99
21) Benzo(a)anthracene	21.266	228	12457	0.410	ng/ul	99
22) Chrysene	21.319	228	14649	0.403	ng/ul	100
24) Benzo(b)fluoranthene	22.850	252	14046	0.424	ng/ul	97
25) Benzo(k)fluoranthene	22.894	252	14710	0.393	ng/ul	99
26) Benzo(a)pyrene	23.426	252	12004	0.431	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.770	276	15493	0.461	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.783	278	12377	0.469	ng/ul	97
29) Benzo(g,h,i)perylene	26.454	276	13240	0.425	ng/ul	99

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

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