

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015428.D
 Acq On : 09 Jul 2021 03:40
 Operator : CG/JU
 Sample : PB137381BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS381

Quant Time: Jul 09 04:35:36 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	7.711	152	2847	0.400	ng/ul	0.00
4) Naphthalene-d8	10.480	136	9438	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.335	164	4900	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	9957	0.400	ng/ul	0.00
17) Chrysene-d12	21.281	240	8138	0.400	ng/ul	0.00
23) Perylene-d12	23.520	264	7102	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	2478	0.772	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	5425	0.366	ng/ul	0.00
18) Fluoranthene-d10	19.118	212	11227	0.420	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.311	88	6459	1.863	ng/ul#	88
5) Naphthalene	10.529	128	11020	0.366	ng/ul	100
7) 2-Methylnaphthalene	12.146	142	7185	0.365	ng/ul	100
8) 1-Methylnaphthalene	12.366	142	6700	0.352	ng/ul	99
10) Acenaphthylene	14.053	152	8661	0.342	ng/ul	99
11) Acenaphthene	14.400	153	7110	0.366	ng/ul	99
12) Fluorene	15.390	166	7875	0.367	ng/ul	98
14) Pentachlorophenol	16.740	266	1619	0.735	ng/ul	99
15) Phenanthrene	17.124	178	13549	0.375	ng/ul	99
16) Anthracene	17.217	178	11356	0.356	ng/ul	100
19) Fluoranthene	19.150	202	15115	0.402	ng/ul	100
20) Pyrene	19.513	202	15337	0.407	ng/ul	100
21) Benzo(a)anthracene	21.267	228	13095	0.434	ng/ul	100
22) Chrysene	21.317	228	14969	0.415	ng/ul	99
24) Benzo(b)fluoranthene	22.848	252	14956	0.451	ng/ul	97
25) Benzo(k)fluoranthene	22.892	252	15081	0.403	ng/ul	98
26) Benzo(a)pyrene	23.421	252	12553	0.450	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.767	276	16215	0.482	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.780	278	12919	0.490	ng/ul	99
29) Benzo(g,h,i)perylene	26.451	276	13956	0.448	ng/ul	98

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

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