

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071221\
 Data File : BN015486.D
 Acq On : 12 Jul 2021 12:53
 Operator : CG/JU
 Sample : PB137440BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS440

Quant Time: Jul 12 13:22:37 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 : Mon Jul 12 11:02:52 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.812	152	3117	0.400	ng/ul	0.00
4) Naphthalene-d8	10.600	136	10284	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.446	164	5139	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.195	188	10015	0.400	ng/ul	0.00
17) Chrysene-d12	21.389	240	7668	0.400	ng/ul	0.00
23) Perylene-d12	23.745	264	6604	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	2414	0.687	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	5017	0.311	ng/ul	0.00
18) Fluoranthene-d10	19.229	212	9334	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	6456	1.700	ng/ul	87
5) Naphthalene	10.650	128	10283	0.313	ng/ul	100
7) 2-Methylnaphthalene	12.267	142	6557	0.305	ng/ul	100
8) 1-Methylnaphthalene	12.481	142	6117	0.295	ng/ul	99
10) Acenaphthylene	14.164	152	7880	0.297	ng/ul	100
11) Acenaphthene	14.506	153	6415	0.315	ng/ul	98
12) Fluorene	15.496	166	6899	0.307	ng/ul	100
14) Pentachlorophenol	16.849	266	1680	0.758	ng/ul	99
15) Phenanthrene	17.233	178	11820	0.325	ng/ul	99
16) Anthracene	17.326	178	9819	0.306	ng/ul	100
19) Fluoranthene	19.257	202	12629	0.356	ng/ul	99
20) Pyrene	19.620	202	12667	0.356	ng/ul	99
21) Benzo(a)anthracene	21.375	228	9961	0.351	ng/ul	99
22) Chrysene	21.427	228	11729	0.345	ng/ul	100
24) Benzo(b)fluoranthene	23.032	252	10709	0.347	ng/ul	98
25) Benzo(k)fluoranthene	23.079	252	11411	0.328	ng/ul	99
26) Benzo(a)pyrene	23.643	252	8871	0.342	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.169	276	12031	0.384	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.192	278	9606	0.391	ng/ul	98
29) Benzo(g,h,i)perylene	26.910	276	9962	0.344	ng/ul	99

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

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