

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010322\
 Data File : BN018316.D
 Acq On : 03 Jan 2022 20:16
 Operator : CG/JU
 Sample : PB141475BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS475

Quant Time: Jan 04 00:06:02 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 28 00:34:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.615	152	1780	0.400	ng/ul	0.00
4) Naphthalene-d8	10.392	136	6715	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.248	164	4975	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.994	188	12071	0.400	ng/ul	0.00
17) Chrysene-d12	21.184	240	18479	0.400	ng/ul	0.00
23) Perylene-d12	23.408	264	20321	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.067	96	1285	0.773	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	3415	0.293	ng/ul	0.00
18) Fluoranthene-d10	19.026	212	15886	0.290	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.105	88	3129	1.692	ng/ul#	83
5) Naphthalene	10.442	128	6407	0.322	ng/ul	97
7) 2-Methylnaphthalene	12.070	142	4292	0.294	ng/ul	100
8) 1-Methylnaphthalene	12.284	142	4527	0.308	ng/ul	98
10) Acenaphthylene	13.966	152	6682	0.311	ng/ul	96
11) Acenaphthene	14.313	153	5987	0.346	ng/ul	98
12) Fluorene	15.303	166	7243	0.339	ng/ul	98
14) Pentachlorophenol	16.656	266	2366	0.495	ng/ul	98
15) Phenanthrene	17.032	178	14045	0.352	ng/ul	99
16) Anthracene	17.129	178	11256	0.314	ng/ul	97
19) Fluoranthene	19.054	202	21530	0.293	ng/ul	98
20) Pyrene	19.416	202	23883	0.311	ng/ul	99
21) Benzo(a)anthracene	21.170	228	23310	0.331	ng/ul	99
22) Chrysene	21.219	228	29721	0.389	ng/ul	100
24) Benzo(b)fluoranthene	22.739	252	30842	0.363	ng/ul	99
25) Benzo(k)fluoranthene	22.783	252	38801	0.400	ng/ul	98
26) Benzo(a)pyrene	23.309	252	30161	0.374	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.662	276	42313	0.412	ng/ul	100
28) Dibenzo(a,h)anthracene	25.682	278	32340	0.408	ng/ul	97
29) Benzo(g,h,i)perylene	26.350	276	37778	0.432	ng/ul	99

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

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