

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041522\
 Data File : BN019469.D
 Acq On : 15 Apr 2022 19:01
 Operator : CG/JU
 Sample : PB144067BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS067

Quant Time: Apr 18 05:04:09 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN041322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 14:35:39 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	3929	0.400	ng/ul	0.00
4) Naphthalene-d8	10.624	136	12501	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.461	164	7608	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.201	188	13758	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	8552	0.400	ng/ul	0.00
23) Perylene-d12	23.748	264	6515	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	3230	0.639	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.219	152	6288	0.356	ng/ul	0.00
18) Fluoranthene-d10	19.235	212	13896	0.439	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.211	88	7754	1.592	ng/ul#	74
5) Naphthalene	10.673	128	12137	0.341	ng/ul	99
7) 2-Methylnaphthalene	12.290	142	7277	0.314	ng/ul#	100
8) 1-Methylnaphthalene	12.505	142	7819	0.322	ng/ul#	100
10) Acenaphthylene	14.179	152	10377	0.376	ng/ul	99
11) Acenaphthene	14.521	153	9007	0.327	ng/ul	98
12) Fluorene	15.511	166	9543	0.311	ng/ul	99
14) Pentachlorophenol	16.868	266	1434	0.649	ng/ul	92
15) Phenanthrene	17.243	178	14978	0.328	ng/ul	100
16) Anthracene	17.340	178	11209	0.319	ng/ul	99
19) Fluoranthene	19.263	202	16633	0.391	ng/ul	99
20) Pyrene	19.626	202	17553	0.388	ng/ul	99
21) Benzo(a)anthracene	21.380	228	9302	0.353	ng/ul	99
22) Chrysene	21.433	228	13129	0.381	ng/ul	99
24) Benzo(b)fluoranthene	23.032	252	10439	0.346	ng/ul	95
25) Benzo(k)fluoranthene	23.078	252	12546	0.360	ng/ul	97
26) Benzo(a)pyrene	23.643	252	9946	0.386	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.176	276	9763	0.341	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.192	278	7147	0.350	ng/ul	91
29) Benzo(g,h,i)perylene	26.917	276	9890	0.357	ng/ul	99

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

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