

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072121\
 Data File : BN015635.D
 Acq On : 20 Jul 2021 23:51
 Operator : CG/JU
 Sample : PB137620BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS620

Quant Time: Jul 21 01:06:41 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 : Tue Jul 20 18:52:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.779	152	2177	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.567	136	7373	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.418	164	3855	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.166	188	8004	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.369	240	6658	0.400	ng/ul	# 0.00
23) Perylene-d12	23.707	264	5832	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.277	96	1961	0.799	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.162	152	4204	0.363	ng/ul	0.00
18) Fluoranthene-d10	19.206	212	9114	0.417	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.306	88	5337	2.013	ng/ul#	71
5) Naphthalene	10.611	128	8524	0.362	ng/ul#	91
7) 2-Methylnaphthalene	12.234	142	5401	0.351	ng/ul	99
8) 1-Methylnaphthalene	12.454	142	5157	0.346	ng/ul	99
10) Acenaphthylene	14.136	152	6724	0.337	ng/ul#	90
11) Acenaphthene	14.478	153	5583	0.366	ng/ul	98
12) Fluorene	15.468	166	6252	0.370	ng/ul	98
14) Pentachlorophenol	16.824	266	781	0.441	ng/ul	92
15) Phenanthrene	17.208	178	10888	0.375	ng/ul	98
16) Anthracene	17.301	178	9051	0.353	ng/ul	95
19) Fluoranthene	19.234	202	11959	0.389	ng/ul#	89
20) Pyrene	19.596	202	12305	0.399	ng/ul#	88
21) Benzo(a)anthracene	21.351	228	9866	0.400	ng/ul	99
22) Chrysene	21.404	228	11358	0.385	ng/ul	98
24) Benzo(b)fluoranthene	22.999	252	10808	0.397	ng/ul	94
25) Benzo(k)fluoranthene	23.046	252	11587	0.377	ng/ul#	93
26) Benzo(a)pyrene	23.605	252	9030	0.395	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.115	276	12390	0.448	ng/ul#	53
28) Dibenzo(a,h)anthracene	26.132	278	9886	0.456	ng/ul#	89
29) Benzo(g,h,i)perylene	26.850	276	10310	0.403	ng/ul#	85

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

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