

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071721\
 Data File : BN015598.D
 Acq On : 16 Jul 2021 15:43
 Operator : CG/JU
 Sample : PB137518BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS518

Quant Time: Jul 16 16:21:10 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 : Fri Jul 16 15:41:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.791	152	2020	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.579	136	6949	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.428	164	3745	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.175	188	7666	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.378	240	6204	0.400	ng/ul	# 0.00
23) Perylene-d12	23.725	264	5402	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.294	96	1989	0.874	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	4492	0.412	ng/ul	0.00
18) Fluoranthene-d10	19.211	212	9235	0.453	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.323	88	5290	2.150	ng/ul#	72
5) Naphthalene	10.628	128	8733	0.394	ng/ul#	91
7) 2-Methylnaphthalene	12.245	142	5661	0.390	ng/ul	99
8) 1-Methylnaphthalene	12.465	142	5425	0.387	ng/ul	99
10) Acenaphthylene	14.145	152	7030	0.363	ng/ul#	91
11) Acenaphthene	14.492	153	5822	0.392	ng/ul	99
12) Fluorene	15.478	166	6502	0.397	ng/ul	98
14) Pentachlorophenol	16.833	266	998	0.588	ng/ul	92
15) Phenanthrene	17.217	178	11041	0.397	ng/ul	98
16) Anthracene	17.310	178	9032	0.368	ng/ul	96
19) Fluoranthene	19.239	202	11962	0.417	ng/ul#	88
20) Pyrene	19.606	202	12113	0.421	ng/ul#	90
21) Benzo(a)anthracene	21.360	228	9476	0.412	ng/ul	99
22) Chrysene	21.413	228	11330	0.412	ng/ul	98
24) Benzo(b)fluoranthene	23.014	252	10723	0.425	ng/ul	94
25) Benzo(k)fluoranthene	23.058	252	11335	0.398	ng/ul#	93
26) Benzo(a)pyrene	23.620	252	8691	0.410	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.136	276	12160	0.475	ng/ul#	52
28) Dibenzo(a,h)anthracene	26.153	278	9588	0.478	ng/ul#	89
29) Benzo(g,h,i)perylene	26.874	276	9899	0.417	ng/ul#	86

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

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