

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016150.D
 Acq On : 02 Sep 2021 23:35
 Operator : CG/JU
 Sample : PB138905BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS905

Quant Time: Sep 03 06:10:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 02 13:32:21 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	2315	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	10028	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.507	164	6016	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.256	188	12073	0.400	ng/ul	0.00
17) Chrysene-d12	21.441	240	12404	0.400	ng/ul	0.00
23) Perylene-d12	23.835	264	12312	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	2067	0.767	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.261	152	6651	0.471	ng/ul	0.00
18) Fluoranthene-d10	19.286	212	16560	0.498	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	5226	1.775	ng/ul#	89
5) Naphthalene	10.716	128	11000	0.412	ng/ul	99
7) 2-Methylnaphthalene	12.333	142	7652	0.439	ng/ul	100
8) 1-Methylnaphthalene	12.547	142	7163	0.412	ng/ul	100
10) Acenaphthylene	14.225	152	11040	0.485	ng/ul	100
11) Acenaphthene	14.567	153	7806	0.410	ng/ul	99
12) Fluorene	15.557	166	8929	0.410	ng/ul	98
14) Pentachlorophenol	16.909	266	2127	1.088	ng/ul	100
15) Phenanthrene	17.294	178	14636	0.421	ng/ul	100
16) Anthracene	17.387	178	14127	0.470	ng/ul	99
19) Fluoranthene	19.314	202	18227	0.441	ng/ul	99
20) Pyrene	19.676	202	18830	0.450	ng/ul	96
21) Benzo(a)anthracene	21.424	228	19281	0.508	ng/ul	98
22) Chrysene	21.479	228	19155	0.447	ng/ul	98
24) Benzo(b)fluoranthene	23.107	252	19676	0.425	ng/ul	93
25) Benzo(k)fluoranthene	23.154	252	19210	0.412	ng/ul#	94
26) Benzo(a)pyrene	23.730	252	17450	0.444	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.316	276	22115	0.436	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.336	278	18505	0.442	ng/ul#	93
29) Benzo(g,h,i)perylene	27.077	276	18703	0.424	ng/ul	95

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

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