

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035341.D
 Acq On : 27 Nov 2024 07:29
 Operator : RC/JU
 Sample : PB165213BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS213

Quant Time: Nov 27 07:56:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 06:09:15 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.309	152	3995	0.400	ng/ul	0.00
4) Naphthalene-d8	10.058	136	10670	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.967	164	6180	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.734	188	11369	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.976	240	7356	0.400	ng/ul	# 0.00
23) Perylene-d12	23.068	264	8190	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.972	96	2248	0.616	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.659	152	5068	0.324	ng/ul	0.00
18) Fluoranthene-d10	18.786	212	7883	0.388	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.001	88	6261	1.580	ng/ul#	87
5) Naphthalene	10.102	128	9003	0.335	ng/ul	99
7) 2-Methylnaphthalene	11.736	142	5980	0.321	ng/ul	98
8) 1-Methylnaphthalene	11.961	142	6085	0.323	ng/ul	98
10) Acenaphthylene	13.680	152	8586	0.350	ng/ul	99
11) Acenaphthene	14.032	153	6512	0.347	ng/ul	99
12) Fluorene	15.031	166	6929	0.315	ng/ul	99
14) Pentachlorophenol	16.392	266	1628	0.615	ng/ul	98
15) Phenanthrene	16.772	178	10218	0.340	ng/ul	99
16) Anthracene	16.865	178	8826	0.316	ng/ul	99
19) Fluoranthene	18.814	202	10034	0.392	ng/ul#	93
20) Pyrene	19.181	202	10087	0.378	ng/ul#	91
21) Benzo(a)anthracene	20.961	228	8590	0.337	ng/ul	100
22) Chrysene	21.011	228	9192	0.358	ng/ul	99
24) Benzo(b)fluoranthene	22.454	252	9380	0.346	ng/ul	96
25) Benzo(k)fluoranthene	22.495	252	10160	0.364	ng/ul#	96
26) Benzo(a)pyrene	22.978	252	8786	0.355	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.108	276	11716	0.351	ng/ul#	86
28) Dibenzo(a,h)anthracene	25.124	278	8867	0.339	ng/ul#	94
29) Benzo(g,h,i)perylene	25.725	276	9363	0.355	ng/ul#	94

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

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