

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035332.D
 Acq On : 27 Nov 2024 01:30
 Operator : RC/JU
 Sample : PB165261BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS261

Quant Time: Nov 27 01:54:01 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	7.310	152	3107	0.400	ng/ul	0.00
4) Naphthalene-d8	10.058	136	8120	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.967	164	4901	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.734	188	9549	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.976	240	5483	0.400	ng/ul	# 0.00
23) Perylene-d12	23.068	264	5907	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.972	96	2102	0.740	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.659	152	4510	0.379	ng/ul	0.00
18) Fluoranthene-d10	18.786	212	7741	0.512	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.001	88	5655	1.834	ng/ul#	88
5) Naphthalene	10.108	128	7798	0.381	ng/ul	100
7) 2-Methylnaphthalene	11.736	142	5319	0.375	ng/ul	99
8) 1-Methylnaphthalene	11.961	142	5368	0.375	ng/ul	99
10) Acenaphthylene	13.680	152	7832	0.403	ng/ul	99
11) Acenaphthene	14.032	153	6047	0.407	ng/ul	97
12) Fluorene	15.036	166	6642	0.380	ng/ul	99
14) Pentachlorophenol	16.396	266	1523	0.685	ng/ul	98
15) Phenanthrene	16.776	178	10065	0.399	ng/ul	100
16) Anthracene	16.869	178	8721	0.372	ng/ul	99
19) Fluoranthene	18.819	202	9767	0.512	ng/ul#	96
20) Pyrene	19.181	202	9654	0.486	ng/ul#	90
21) Benzo(a)anthracene	20.961	228	7354	0.387	ng/ul	98
22) Chrysene	21.011	228	8251	0.431	ng/ul	100
24) Benzo(b)fluoranthene	22.454	252	7778	0.398	ng/ul	96
25) Benzo(k)fluoranthene	22.498	252	8912	0.443	ng/ul#	96
26) Benzo(a)pyrene	22.978	252	7126	0.399	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.111	276	9856	0.409	ng/ul#	86
28) Dibenzo(a,h)anthracene	25.128	278	7493	0.397	ng/ul#	94
29) Benzo(g,h,i)perylene	25.725	276	7927	0.417	ng/ul#	94

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

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