

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035179.D
 Acq On : 20 Nov 2024 01:28
 Operator : RC/JU
 Sample : PB165072BS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS072

Quant Time: Nov 20 01:52:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3356	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	7794	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.185	164	4962	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	10429	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	9486	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	10467	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1986	0.647	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	4171	0.365	ng/ul	0.00
18) Fluoranthene-d10	18.977	212	9663	0.369	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	5396	1.621	ng/ul#	95
5) Naphthalene	10.355	128	6747	0.344	ng/ul	100
7) 2-Methylnaphthalene	11.978	142	4638	0.341	ng/ul	99
8) 1-Methylnaphthalene	12.203	142	4690	0.341	ng/ul	100
10) Acenaphthylene	13.898	152	6765	0.344	ng/ul	99
11) Acenaphthene	14.245	153	5201	0.345	ng/ul	97
12) Fluorene	15.240	166	5993	0.339	ng/ul	99
14) Pentachlorophenol	16.595	266	1667	0.686	ng/ul	98
15) Phenanthrene	16.979	178	9600	0.348	ng/ul	100
16) Anthracene	17.072	178	8720	0.341	ng/ul	99
19) Fluoranthene	19.009	202	11482	0.348	ng/ul	99
20) Pyrene	19.372	202	12078	0.351	ng/ul	100
21) Benzo(a)anthracene	21.134	228	11944	0.363	ng/ul	100
22) Chrysene	21.183	228	12448	0.376	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	12713	0.367	ng/ul	98
25) Benzo(k)fluoranthene	22.714	252	13489	0.378	ng/ul	97
26) Benzo(a)pyrene	23.226	252	11640	0.368	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.477	276	15977	0.374	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.493	278	12535	0.374	ng/ul	98
29) Benzo(g,h,i)perylene	26.131	276	12648	0.376	ng/ul	99

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

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