

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111824\
 Data File : BN035149.D
 Acq On : 18 Nov 2024 15:45
 Operator : RC/JU
 Sample : PB164837BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS837

Quant Time: Nov 18 16:25:54 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	7.545	152	3566	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	8795	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.189	164	5827	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.941	188	12283	0.400	ng/ul	0.00
17) Chrysene-d12	21.151	240	10489	0.400	ng/ul	0.00
23) Perylene-d12	23.325	264	11098	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	2739	0.840	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	6451	0.501	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	14793	0.511	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.170	88	7402	2.092	ng/ul#	94
5) Naphthalene	10.355	128	9828	0.444	ng/ul	99
7) 2-Methylnaphthalene	11.983	142	6846	0.446	ng/ul	99
8) 1-Methylnaphthalene	12.203	142	7937	0.512	ng/ul	99
10) Acenaphthylene	13.902	152	10184	0.441	ng/ul	100
11) Acenaphthene	14.249	153	7837	0.443	ng/ul	98
12) Fluorene	15.244	166	9053	0.436	ng/ul	99
14) Pentachlorophenol	16.595	266	2411	0.843	ng/ul	99
15) Phenanthrene	16.983	178	14469	0.446	ng/ul	99
16) Anthracene	17.072	178	13217	0.439	ng/ul	99
19) Fluoranthene	19.009	202	16859	0.462	ng/ul	100
20) Pyrene	19.376	202	17415	0.458	ng/ul	99
21) Benzo(a)anthracene	21.137	228	16439	0.452	ng/ul	100
22) Chrysene	21.186	228	16917	0.462	ng/ul	99
24) Benzo(b)fluoranthene	22.676	252	16856	0.459	ng/ul	97
25) Benzo(k)fluoranthene	22.720	252	17755	0.469	ng/ul	96
26) Benzo(a)pyrene	23.229	252	15364	0.458	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.483	276	20763	0.459	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.500	278	16256	0.458	ng/ul	97
29) Benzo(g,h,i)perylene	26.137	276	16242	0.455	ng/ul	98

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

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