

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111524\
 Data File : BN035112.D
 Acq On : 15 Nov 2024 18:48
 Operator : RC/JU
 Sample : PB164837BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS837

Quant Time: Nov 15 22:27:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 00:41:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.550	152	2684	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	7030	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.185	164	5657	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	14042	0.400	ng/ul	0.00
17) Chrysene-d12	21.137	240	11950	0.400	ng/ul	0.00
23) Perylene-d12	23.299	264	10179	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	2121	0.830	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	5535	0.514	ng/ul	-0.01
18) Fluoranthene-d10	18.972	212	17891	0.583	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.170	88	5678	2.029	ng/ul#	90
5) Naphthalene	10.361	128	7930	0.451	ng/ul	98
7) 2-Methylnaphthalene	11.983	142	5867	0.447	ng/ul	99
8) 1-Methylnaphthalene	12.203	142	6891	0.523	ng/ul	100
10) Acenaphthylene	13.903	152	9774	0.423	ng/ul	100
11) Acenaphthene	14.249	153	7590	0.433	ng/ul	98
12) Fluorene	15.244	166	9410	0.422	ng/ul	98
14) Pentachlorophenol	16.591	266	2919	0.899	ng/ul	97
15) Phenanthrene	16.975	178	16552	0.441	ng/ul	98
16) Anthracene	17.068	178	15126	0.429	ng/ul	98
19) Fluoranthene	19.000	202	20489	0.513	ng/ul#	96
20) Pyrene	19.362	202	21172	0.511	ng/ul#	95
21) Benzo(a)anthracene	21.119	228	18996	0.439	ng/ul	99
22) Chrysene	21.172	228	19091	0.432	ng/ul	99
24) Benzo(b)fluoranthene	22.656	252	16586	0.463	ng/ul	98
25) Benzo(k)fluoranthene	22.700	252	17192	0.459	ng/ul#	98
26) Benzo(a)pyrene	23.206	252	14443	0.452	ng/ul#	98
27) Indeno(1,2,3-cd)pyrene	25.450	276	16351	0.418	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.463	278	12758	0.419	ng/ul#	98
29) Benzo(g,h,i)perylene	26.104	276	12726	0.426	ng/ul#	97

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

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