

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121421\
 Data File : BM033477.D
 Acq On : 15 Dec 2021 13:57
 Operator : CG/JU
 Sample : PB141312BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS312

Quant Time: Dec 16 02:33:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 15 01:00:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.901	152	573	0.400	ng/ul	0.00
4) Naphthalene-d8	10.702	136	1550	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.523	164	1025	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.268	188	2344	0.400	ng/ul	-0.01
17) Chrysene-d12	21.429	240	2645	0.400	ng/ul	0.00
23) Perylene-d12	23.745	264	2158	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.362	96	374	0.806	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.294	152	796	0.407	ng/ul	0.00
18) Fluoranthene-d10	19.285	212	2861	0.497	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.396	88	1204	1.984	ng/ul#	83
5) Naphthalene	10.752	128	1942	0.342	ng/ul	98
7) 2-Methylnaphthalene	12.366	142	1201	0.372	ng/ul	94
8) 1-Methylnaphthalene	12.573	142	1206	0.361	ng/ul	100
10) Acenaphthylene	14.246	152	1922	0.335	ng/ul	98
11) Acenaphthene	14.583	153	1550	0.344	ng/ul	97
12) Fluorene	15.577	166	1731	0.359	ng/ul	98
14) Pentachlorophenol	16.934	266	439	0.522	ng/ul	99
15) Phenanthrene	17.309	178	2791	0.365	ng/ul	97
16) Anthracene	17.406	178	2490	0.369	ng/ul	98
19) Fluoranthene	19.311	202	3484	0.501	ng/ul#	93
20) Pyrene	19.674	202	3699	0.482	ng/ul#	89
21) Benzo(a)anthracene	21.417	228	2479	0.421	ng/ul	99
22) Chrysene	21.468	228	3078	0.419	ng/ul	99
24) Benzo(b)fluoranthene	23.044	252	2604	0.454	ng/ul	89
25) Benzo(k)fluoranthene	23.088	252	2921	0.443	ng/ul#	90
26) Benzo(a)pyrene	23.650	252	2233	0.404	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.112	276	2689	0.435	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.143	278	2246	0.439	ng/ul	92
29) Benzo(g,h,i)perylene	26.829	276	2437	0.438	ng/ul#	94

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

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