

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112423\
 Data File : BM043025.D
 Acq On : 25 Nov 2023 21:03
 Operator : MA/JU
 Sample : PB157276BS
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS276

Quant Time: Nov 27 02:44:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 27 02:36:19 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.872	152	539	0.400	ng/ul	0.00
4) Naphthalene-d8	10.644	136	1870	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.464	164	1148	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.229	188	2400	0.400	ng/ul	0.00
17) Chrysene-d12	21.412	240	1391	0.400	ng/ul	-0.01
23) Perylene-d12	23.768	264	1742	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	827	1.000	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.244	152	1350	0.540	ng/ul	0.00
18) Fluoranthene-d10	19.256	212	2614	0.516	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.197	88	2332	2.626	ng/ul#	84
5) Naphthalene	10.693	128	2305	0.454	ng/ul	97
7) 2-Methylnaphthalene	12.315	142	1437	0.479	ng/ul	97
8) 1-Methylnaphthalene	12.513	142	1430	0.428	ng/ul	95
10) Acenaphthylene	14.195	152	2644	0.451	ng/ul	97
11) Acenaphthene	14.528	153	2022	0.456	ng/ul	98
12) Fluorene	15.532	166	2138	0.469	ng/ul	98
14) Pentachlorophenol	16.882	266	581	0.894	ng/ul	97
15) Phenanthrene	17.271	178	3051	0.445	ng/ul	99
16) Anthracene	17.381	178	2943	0.454	ng/ul	97
19) Fluoranthene	19.284	202	3692	0.468	ng/ul	98
20) Pyrene	19.651	202	3862	0.461	ng/ul	98
21) Benzo(a)anthracene	21.403	228	2217	0.491	ng/ul	97
22) Chrysene	21.450	228	3744	0.445	ng/ul	99
24) Benzo(b)fluoranthene	23.040	252	2998	0.510	ng/ul	95
25) Benzo(k)fluoranthene	23.095	252	4096	0.506	ng/ul	98
26) Benzo(a)pyrene	23.668	252	3153	0.491	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.229	276	4384	0.505	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.249	278	3455	0.514	ng/ul	93
29) Benzo(g,h,i)perylene	26.997	276	4288	0.508	ng/ul	97

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

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