

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041024\
 Data File : BM045119.D
 Acq On : 11 Apr 2024 03:38
 Operator : MA/JU
 Sample : PB160083BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS083

Quant Time: Apr 11 04:08:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 10 11:09:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.648	152	1201	0.400	ng/ul	0.00
4) Naphthalene-d8	10.424	136	3728	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.280	164	1969	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.052	188	4146	0.400	ng/ul	0.00
17) Chrysene-d12	21.256	240	2806	0.400	ng/ul #	0.00
23) Perylene-d12	23.478	264	3492	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.209	96	1310	0.830	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.041	152	1961	0.388	ng/ul	0.01
18) Fluoranthene-d10	19.086	212	3482	0.491	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.239	88	3185	2.106	ng/ul#	82
5) Naphthalene	10.479	128	3907	0.395	ng/ul	95
7) 2-Methylnaphthalene	12.112	142	2296	0.389	ng/ul	97
8) 1-Methylnaphthalene	12.316	142	2400	0.372	ng/ul	95
10) Acenaphthylene	14.007	152	3857	0.393	ng/ul	96
11) Acenaphthene	14.345	153	2801	0.398	ng/ul	98
12) Fluorene	15.358	166	2966	0.376	ng/ul	97
14) Pentachlorophenol	16.719	266	644	0.662	ng/ul	99
15) Phenanthrene	17.095	178	4363	0.374	ng/ul	98
16) Anthracene	17.200	178	4406	0.383	ng/ul	94
19) Fluoranthene	19.118	202	4948	0.493	ng/ul	96
20) Pyrene	19.481	202	5276	0.498	ng/ul	96
21) Benzo(a)anthracene	21.251	228	3647	0.373	ng/ul	96
22) Chrysene	21.294	228	5338	0.438	ng/ul	97
24) Benzo(b)fluoranthene	22.811	252	4748	0.376	ng/ul	93
25) Benzo(k)fluoranthene	22.858	252	6122	0.433	ng/ul#	88
26) Benzo(a)pyrene	23.390	252	5039	0.418	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.732	276	6386	0.387	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.748	278	4889	0.374	ng/ul#	86
29) Benzo(g,h,i)perylene	26.389	276	5942	0.417	ng/ul	96

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

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