

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041024\
 Data File : BM045093.D
 Acq On : 10 Apr 2024 11:20
 Operator : MA/JU
 Sample : P2013-15MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0MC0MS

Quant Time: Apr 10 11:49:10 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	7.639	152	1279	0.400	ng/ul	0.00
4) Naphthalene-d8	10.418	136	4251	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.279	164	2236	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.043	188	4656	0.400	ng/ul	-0.01
17) Chrysene-d12	21.249	240	3535	0.400	ng/ul	# 0.00
23) Perylene-d12	23.467	264	4168	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	536	0.319	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.019	152	2041	0.354	ng/ul	-0.01
18) Fluoranthene-d10	19.075	212	4810	0.538	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.243	88	1252	0.777	ng/ul#	83
5) Naphthalene	10.468	128	4084	0.362	ng/ul	98
7) 2-Methylnaphthalene	12.090	142	2525	0.375	ng/ul	98
8) 1-Methylnaphthalene	12.305	142	2587	0.352	ng/ul	98
10) Acenaphthylene	14.001	152	4389	0.393	ng/ul	96
11) Acenaphthene	14.339	153	3206	0.401	ng/ul	100
12) Fluorene	15.348	166	3499	0.391	ng/ul	99
14) Pentachlorophenol	16.701	266	1055	0.966	ng/ul	97
15) Phenanthrene	17.081	178	5616	0.428	ng/ul	99
16) Anthracene	17.182	178	5527	0.428	ng/ul	97
19) Fluoranthene	19.108	202	6864	0.543	ng/ul	98
20) Pyrene	19.470	202	7178	0.538	ng/ul	98
21) Benzo(a)anthracene	21.234	228	5268	0.428	ng/ul	98
22) Chrysene	21.284	228	6786	0.442	ng/ul	98
24) Benzo(b)fluoranthene	22.800	252	6643	0.441	ng/ul	94
25) Benzo(k)fluoranthene	22.847	252	7501	0.445	ng/ul#	91
26) Benzo(a)pyrene	23.373	252	6414	0.446	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	25.706	276	8322	0.422	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.732	278	6235	0.400	ng/ul#	90
29) Benzo(g,h,i)perylene	26.380	276	7526	0.443	ng/ul	96

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

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