

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103124\
 Data File : BM048415.D
 Acq On : 02 Nov 2024 03:47
 Operator : RC/JU
 Sample : P4558-08MS
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E1L92MS

Quant Time: Nov 03 21:52:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.788	152	2382	0.400	ng/ul	0.04
4) Naphthalene-d8	10.551	136	9484	0.400	ng/ul	0.02
9) Acenaphthene-d10	14.363	164	5637	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.120	188	10595	0.400	ng/ul	0.00
17) Chrysene-d12	21.289	240	10171	0.400	ng/ul	0.00
23) Perylene-d12	23.528	264	10160	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	1241	0.424	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.152	152	4638	0.384	ng/ul	0.01
18) Fluoranthene-d10	19.137	212	13909	0.539	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.197	88	3763	1.186	ng/ul#	89
5) Naphthalene	10.595	128	8809	0.355	ng/ul	96
7) 2-Methylnaphthalene	12.229	142	4847	0.329	ng/ul	97
8) 1-Methylnaphthalene	12.421	142	5898	0.375	ng/ul	98
10) Acenaphthylene	14.095	152	7981	0.366	ng/ul	98
11) Acenaphthene	14.428	153	6749	0.386	ng/ul	98
12) Fluorene	15.436	166	7521	0.387	ng/ul	97
14) Pentachlorophenol	16.782	266	2643	0.896	ng/ul	95
15) Phenanthrene	17.162	178	12045	0.457	ng/ul	99
16) Anthracene	17.268	178	10543	0.428	ng/ul	98
19) Fluoranthene	19.169	202	17472	0.498	ng/ul	99
20) Pyrene	19.527	202	18538	0.493	ng/ul	100
21) Benzo(a)anthracene	21.280	228	12690	0.417	ng/ul	99
22) Chrysene	21.324	228	21572	0.514	ng/ul	100
24) Benzo(b)fluoranthene	22.852	252	16248	0.468	ng/ul	98
25) Benzo(k)fluoranthene	22.896	252	21378	0.520	ng/ul	100
26) Benzo(a)pyrene	23.437	252	15207	0.466	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.816	276	21051	0.462	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.836	278	16209	0.456	ng/ul	100
29) Benzo(g,h,i)perylene	26.496	276	18410	0.479	ng/ul	98

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

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