

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110722\
 Data File : BM037409.D
 Acq On : 08 Nov 2022 02:14
 Operator : CG/JU
 Sample : N5408-14MSD
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4Z2MSD

Quant Time: Nov 08 03:21:24 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 00:30:18 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.837	152	4478	0.400	ng/ul	0.00
4) Naphthalene-d8	10.633	136	14493	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.468	164	8424	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.208	188	17497	0.400	ng/ul	0.00
17) Chrysene-d12	21.380	240	11836	0.400	ng/ul	0.00
23) Perylene-d12	23.715	264	9906	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.259	96	1169	0.206	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.222	152	6554	0.301	ng/ul	0.00
18) Fluoranthene-d10	19.233	212	17623	0.476	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	3199	0.569	ng/ul#	90
5) Naphthalene	10.682	128	13161	0.313	ng/ul	98
7) 2-Methylnaphthalene	12.299	142	8098	0.308	ng/ul	96
8) 1-Methylnaphthalene	12.514	142	8795	0.325	ng/ul	99
10) Acenaphthylene	14.191	152	11678	0.326	ng/ul	100
11) Acenaphthene	14.529	153	10411	0.328	ng/ul	98
12) Fluorene	15.519	166	12257	0.338	ng/ul	99
14) Pentachlorophenol	16.866	266	3471	0.752	ng/ul	99
15) Phenanthrene	17.250	178	22120	0.369	ng/ul	100
16) Anthracene	17.343	178	18321	0.375	ng/ul	99
19) Fluoranthene	19.261	202	27363	0.488	ng/ul	97
20) Pyrene	19.624	202	28644	0.488	ng/ul	100
21) Benzo(a)anthracene	21.365	228	22202	0.435	ng/ul	100
22) Chrysene	21.415	228	24378	0.428	ng/ul	100
24) Benzo(b)fluoranthene	23.002	252	24836	0.437	ng/ul	97
25) Benzo(k)fluoranthene	23.049	252	23516	0.432	ng/ul	98
26) Benzo(a)pyrene	23.610	252	19738	0.415	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.121	276	25113	0.397	ng/ul	98
28) Dibenzo(a,h)anthracene	26.142	278	19400	0.398	ng/ul	97
29) Benzo(g,h,i)perylene	26.863	276	21659	0.396	ng/ul	98

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

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