

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121522\
 Data File : BM038061.D
 Acq On : 16 Dec 2022 10:30
 Operator : CG/JU
 Sample : N5925-01
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBXP3

Quant Time: Dec 16 11:13:23 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 15 21:53:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	4004	0.400	ng/ul	0.00
4) Naphthalene-d8	10.873	136	11863	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.680	164	6185	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.417	188	13927	0.400	ng/ul	0.00
17) Chrysene-d12	21.586	240	11762	0.400	ng/ul	0.00
23) Perylene-d12	24.038	264	12598	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	2263	0.414	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.451	152	4354	0.251	ng/ul	0.00
18) Fluoranthene-d10	19.436	212	12875	0.302	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.922	128	7443	0.207	ng/ul	99
7) 2-Methylnaphthalene	12.522	142	9852	0.446	ng/ul	99
10) Acenaphthylene	14.402	152	7173	0.221	ng/ul	97
11) Acenaphthene	14.745	153	7011	0.285	ng/ul	98
12) Fluorene	15.726	166	11289	0.415	ng/ul	99
14) Pentachlorophenol	17.067	266	4015	0.696	ng/ul	100
15) Phenanthrene	17.460	178	15359	0.327	ng/ul	99
16) Anthracene	17.552	178	18701	0.431	ng/ul	99
19) Fluoranthene	19.464	202	20242	0.333	ng/ul	98
20) Pyrene	19.826	202	20425	0.329	ng/ul	99
21) Benzo(a)anthracene	21.568	228	24791	0.488	ng/ul	99
22) Chrysene	21.624	228	23873	0.466	ng/ul	99
24) Benzo(b)fluoranthene	23.284	252	32666	0.547	ng/ul	96
25) Benzo(k)fluoranthene	23.336	252	16600	0.288	ng/ul	98
26) Benzo(a)pyrene	23.927	252	20487	0.407	ng/ul	97
28) Dibenzo(a,h)anthracene	26.622	278	33379	0.678	ng/ul	97
29) Benzo(g,h,i)perylene	27.397	276	14516	0.271	ng/ul	98

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

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