

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090122\
 Data File : BM036479.D
 Acq On : 01 Sep 2022 18:49
 Operator : CG/JU
 Sample : N4458-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0KW3

Quant Time: Sep 02 01:37:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 02 01:33:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	179592	20.000	ng/u1	0.00
20) Naphthalene-d8	10.522	136	842680	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.374	164	531180	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.121	188	1113950	20.000	ng/u1	0.00
79) Chrysene-d12	21.309	240	1206181	20.000	ng/u1	0.00
88) Perylene-d12	23.603	264	1183231	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	21920	4.652	ng/uL	0.00
4) Pyridine-d5	3.563	84	111706	8.596	ng/u1	0.00
7) Phenol-d5	6.904	99	131249	8.562	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	348375	36.104	ng/u1	0.00
11) 2-Chlorophenol-d4	7.257	132	324696	26.958	ng/u1	0.00
15) 4-Methylphenol-d8	8.451	113	275817	24.097	ng/u1	0.00
21) Nitrobenzene-d5	8.898	128	251081	39.298	ng/u1	0.00
24) 2-Nitrophenol-d4	9.616	143	254608	39.287	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.151	165	407445	35.186	ng/u1	0.00
31) 4-Chloroaniline-d4	10.681	131	206361	11.495	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	1539562	45.814	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	1820223	43.428	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	35597	5.590	ng/u1	0.01
60) Fluorene-d10	15.369	176	1396061	45.974	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	226484	39.547	ng/u1	0.00
73) Anthracene-d10	17.221	188	2395673	48.367	ng/u1	0.00
81) Pyrene-d10	19.521	212	2970810	47.480	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.456	264	2872301	49.511	ng/u1	0.00

Target Compounds	Qvalue

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

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