

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036569.D
 Acq On : 16 Sep 2022 18:58
 Operator : CG/JU
 Sample : N4601-17
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5758

Quant Time: Sep 16 23:02:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	334355	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	1479275	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	978408	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	2100734	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	2087945	20.000	ng/u1	0.00
88) Perylene-d12	23.997	264	2168443	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	21987	2.465	ng/uL	0.00
4) Pyridine-d5	3.822	84	65215	2.507	ng/u1	0.00
7) Phenol-d5	7.175	99	370797	12.410	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	277418	15.288	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	333615	15.508	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	336733	15.252	ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	167819	16.761	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	164037	18.291	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	335104	16.123	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	460214	13.318	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	1435916	21.760	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	1172201	14.976	ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	194500	15.197	ng/u1	0.00
60) Fluorene-d10	15.639	176	1140733	19.290	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	111856	13.458	ng/u1	0.00
73) Anthracene-d10	17.486	188	1653969	17.220	ng/u1	0.00
81) Pyrene-d10	19.762	212	2239092	21.798	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.838	264	2334946	22.040	ng/u1	0.00

Target Compounds	Qvalue

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

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