

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036668.D
 Acq On : 22 Sep 2022 19:05
 Operator : CG/JU
 Sample : SP6011
 Misc : SFAM SURROGATE
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP6011

Quant Time: Sep 23 01:04:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 21 22:42:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	459253	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1975957	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1209915	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2347061	20.000	ng/u1	0.00
79) Chrysene-d12	21.544	240	1775121	20.000	ng/u1	0.00
88) Perylene-d12	23.985	264	1566713	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	157355	12.844	ng/uL	0.00
4) Pyridine-d5	3.816	84	2360017	66.054	ng/u1	0.00
7) Phenol-d5	7.181	99	2959533	72.112	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	1779991	71.417	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	2307498	78.094	ng/u1	0.00
15) 4-Methylphenol-d8	8.734	113	2314870	76.335	ng/u1	0.00
21) Nitrobenzene-d5	9.186	128	1180307	88.252	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	1233584	102.973	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	2153737	77.577	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	3307427	71.653	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	6337941	77.669	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	7564436	78.150	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	1115043	70.450	ng/u1	0.01
60) Fluorene-d10	15.633	176	5623973	76.906	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	742128	79.919	ng/u1	0.00
73) Anthracene-d10	17.480	188	8214867	76.550	ng/u1	0.00
81) Pyrene-d10	19.762	212	8132313	93.123	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.838	264	6037655	78.879	ng/u1	0.00

Target Compounds Qvalue

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

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