

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112822\
 Data File : BM037782.D
 Acq On : 29 Nov 2022 06:06
 Operator : CG/JU
 Sample : SP6063
 Misc : SFAM SURROGATE
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP6063

Quant Time: Nov 29 06:53:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 28 22:24:47 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/29/2022
 Supervised By :mohammad ahmed 11/30/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.104	152	180939	20.000	ng/u1	0.00
20) Naphthalene-d8	10.927	136	824784	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.727	164	482178	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.462	188	929295	20.000	ng/u1	0.00
79) Chrysene-d12	21.633	240	771682	20.000	ng/u1	0.00
88) Perylene-d12	24.121	264	619839	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	72759	16.676	ng/uL	0.00
4) Pyridine-d5	3.863	84	1216054	85.629	ng/u1	0.00
7) Phenol-d5	7.257	99	1485058	82.268	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	1040018	86.368	ng/u1	0.00
11) 2-Chlorophenol-d4	7.634	132	1157535	92.283	ng/u1	0.00
15) 4-Methylphenol-d8	8.816	113	1172012	81.590	ng/u1	0.00
21) Nitrobenzene-d5	9.281	128	622975	96.348	ng/u1	0.00
24) 2-Nitrophenol-d4	10.004	143	650318	102.072	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.539	165	1044151	89.555	ng/u1	0.00
31) 4-Chloroaniline-d4	11.063	131	1654723	85.527	ng/u1	0.00
46) Dimethylphthalate-d6	14.139	166	3267290	95.664	ng/u1	0.00
49) Acenaphthylene-d8	14.427	160	3920354	97.248	ng/u1	0.00
54) 4-Nitrophenol-d4	14.915	143	617299	82.055	ng/u1	0.00
60) Fluorene-d10	15.721	176	2851299	95.323	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	442591	88.535	ng/u1	0.00
73) Anthracene-d10	17.568	188	4394776m	101.129	ng/u1	0.00
81) Pyrene-d10	19.850	212	4853682	123.771	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.962	264	3168333	99.306	ng/u1	0.00

Target Compounds Qvalue

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

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