

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
 Data File : BM030874.D
 Acq On : 08 Jul 2021 02:26
 Operator : CG/JU
 Sample : SP5610
 Misc : SFAM SURROGATE
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP5610

Quant Time: Jul 08 06:05:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 08 03:09:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	81231	20.000	ng/u1	0.00
20) Naphthalene-d8	10.545	136	338861	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.392	164	202035	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.127	188	363612	20.000	ng/u1	-0.01
79) Chrysene-d12	21.303	240	290640	20.000	ng/u1	0.00
88) Perylene-d12	23.544	264	315264	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	34970	16.822	ng/uL	-0.01
4) Pyridine-d5	3.663	84	486621	89.514	ng/u1	-0.01
7) Phenol-d5	6.934	99	571985	80.586	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.098	67	382365	83.633	ng/u1	0.00
11) 2-Chlorophenol-d4	7.292	132	465345	85.453	ng/u1	-0.01
15) 4-Methylphenol-d8	8.463	113	463442	81.475	ng/u1	-0.01
21) Nitrobenzene-d5	8.916	128	225079	89.306	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.633	143	246455	89.372	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.169	165	421783	79.479	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.686	131	621171	82.424	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	1222485	86.430	ng/u1	0.00
49) Acenaphthylene-d8	14.086	160	1612808	90.269	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	167573	68.077	ng/u1	0.00
60) Fluorene-d10	15.386	176	1076211	86.410	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	140574	60.292	ng/u1	0.00
73) Anthracene-d10	17.227	188	1529492	90.266	ng/u1	-0.01
81) Pyrene-d10	19.521	212	1504649	93.868	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.403	264	1438788	88.413	ng/u1	-0.01

Target Compounds Qvalue

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

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