

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037364.D
 Acq On : 04 Nov 2022 02:09
 Operator : CG/JU
 Sample : PB148654BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK654

Quant Time: Nov 04 02:47:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:39:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	307994	20.000	ng/u1	0.00
20) Naphthalene-d8	10.645	136	1385245	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.480	164	908161	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	1855420	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1663308	20.000	ng/u1	0.00
88) Perylene-d12	23.738	264	1434801	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	54530	7.805	ng/uL	0.00
4) Pyridine-d5	3.693	84	757312	36.310	ng/u1	0.00
7) Phenol-d5	7.022	99	954754	35.620	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	609755	37.072	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	790056	38.578	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	770071	35.183	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	389561	37.647	ng/u1	0.00
24) 2-Nitrophenol-d4	9.734	143	421327	37.380	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	738574	35.541	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	1113431	35.062	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	2300334	36.634	ng/u1	0.00
49) Acenaphthylene-d8	14.174	160	2808297	37.201	ng/u1	0.00
54) 4-Nitrophenol-d4	14.698	143	382638	30.486	ng/u1	0.00
60) Fluorene-d10	15.474	176	2190217	39.114	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	328212	27.929	ng/u1	0.00
73) Anthracene-d10	17.321	188	3414755	39.962	ng/u1	0.00
81) Pyrene-d10	19.609	212	3789201	43.595	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.586	264	3061596	41.453	ng/u1	0.00

Target Compounds Qvalue

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

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