

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036671.D
 Acq On : 23 Sep 2022 11:25
 Operator : CG/JU
 Sample : PB147811BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK811

Quant Time: Sep 23 23:34:37 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	460906	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2090409	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	1322753	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.374	188	2484035	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	1733636	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	1543915	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	60552	4.925	ng/uL	0.00
4) Pyridine-d5	3.816	84	771683	21.521	ng/ul	0.00
7) Phenol-d5	7.169	99	1084706	26.335	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	673806	26.938	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	869337	29.316	ng/ul	-0.01
15) 4-Methylphenol-d8	8.722	113	875988	28.783	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	433893	30.666	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	429321	33.875	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.439	165	801680	27.295	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.963	131	1210956	24.798	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	2520718	28.255	ng/ul	0.00
49) Acenaphthylene-d8	14.333	160	2934488	27.731	ng/ul	0.00
54) 4-Nitrophenol-d4	14.827	143	370403	21.406	ng/ul	0.00
60) Fluorene-d10	15.627	176	2183030	27.306	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	210811	21.450	ng/ul	0.00
73) Anthracene-d10	17.474	188	3235589	28.488	ng/ul	0.00
81) Pyrene-d10	19.756	212	3167292	37.136	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	2286486	30.313	ng/ul	0.00

Target Compounds Qvalue

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

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