

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036648.D
 Acq On : 22 Sep 2022 05:33
 Operator : CG/JU
 Sample : N4649-08
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ36

Quant Time: Sep 22 06:17:31 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	8.022	152	409024	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	1850944	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1161840	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2286214	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	1665660	20.000	ng/u1	0.00
88) Perylene-d12	23.986	264	1451874	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	43197	3.959	ng/uL	0.00
4) Pyridine-d5	3.828	84	216280	6.797	ng/u1	0.01
7) Phenol-d5	7.175	99	247419	6.769	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	686450	30.924	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	699201	26.570	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	468654	17.352	ng/u1	0.00
21) Nitrobenzene-d5	9.187	128	464639	37.088	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	449716	40.075	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	785123	30.190	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1074808	24.858	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2831948	36.140	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	3194861	34.373	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	91663	6.031	ng/u1	0.00
60) Fluorene-d10	15.633	176	2514490	35.808	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	310800	34.361	ng/u1	0.00
73) Anthracene-d10	17.480	188	4006966	38.333	ng/u1	0.00
81) Pyrene-d10	19.762	212	4160047	50.767	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2839933	40.037	ng/u1	0.00

Target Compounds Qvalue

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

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