

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036729.D
 Acq On : 25 Sep 2022 00:47
 Operator : CG/JU
 Sample : N4662-07
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ28

Quant Time: Sep 26 01:15:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 24 00:41:16 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	364410	20.000	ng/u1	0.00
20) Naphthalene-d8	10.816	136	1633292	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.627	164	1034764	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.368	188	2076276	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1534699	20.000	ng/u1	-0.01
88) Perylene-d12	23.962	264	1346958	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	36525	3.757	ng/uL	0.00
4) Pyridine-d5	3.822	84	110603	3.901	ng/u1	0.00
7) Phenol-d5	7.157	99	183903	5.647	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.328	67	539140	27.261	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.534	132	528643	22.548	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	351814	14.621	ng/u1	-0.02
21) Nitrobenzene-d5	9.169	128	346506	31.344	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.892	143	317034	32.017	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.434	165	593659	25.870	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.951	131	569588	14.929	ng/u1	-0.01
46) Dimethylphthalate-d6	14.039	166	2138701	30.645	ng/u1	-0.01
49) Acenaphthylene-d8	14.322	160	2486008	30.031	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.816	143	56010	4.138	ng/u1	-0.02
60) Fluorene-d10	15.616	176	1987621	31.781	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.727	200	202946	24.705	ng/u1	-0.01
73) Anthracene-d10	17.462	188	3240784	34.138	ng/u1	-0.01
81) Pyrene-d10	19.751	212	3434744	45.493	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.803	264	2337986	35.528	ng/u1	-0.02

Target Compounds Qvalue

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

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