

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
 Data File : BP010787.D
 Acq On : 24 Jun 2022 11:38
 Operator : CG/JU
 Sample : N3395-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0157

Quant Time: Jun 24 22:44:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	360994	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	1465349	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.375	164	780968	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.139	188	1404854	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.251	240	1168450	20.000	ng/u1	# 0.00
88) Perylene-d12	23.627	264	1306355	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	57536	6.029	ng/uL	0.00
4) Pyridine-d5	3.611	84	331604	13.054	ng/u1	0.00
7) Phenol-d5	6.887	99	368537	12.896	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	760577	40.697	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	771136	32.525	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	533530	23.618	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	421021	45.133	ng/u1	0.00
24) 2-Nitrophenol-d4	9.604	143	345734	44.798	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.140	165	704940	36.968	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1052104	33.800	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	2352788	43.980	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2878654	42.162	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	77187	9.237	ng/u1	0.00
60) Fluorene-d10	15.381	176	1979153	43.111	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	150933	32.428	ng/u1	0.00
73) Anthracene-d10	17.239	188	2949878	47.195	ng/u1	0.00
81) Pyrene-d10	19.492	212	3048812	47.894	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.474	264	3108808	48.158	ng/u1	0.00

Target Compounds Qvalue

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

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