

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
 Data File : BP010913.D
 Acq On : 30 Jun 2022 10:56
 Operator : CG/JU
 Sample : N3496-10
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 01 02:32:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 27 23:12:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	538551	20.000	ng/u1	0.00
20) Naphthalene-d8	10.498	136	2203115	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	1126580	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.127	188	2037202	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1703332	20.000	ng/u1	# 0.00
88) Perylene-d12	23.598	264	1797929	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	310	0.022	ng/uL	0.00
4) Pyridine-d5	3.616	84	641	0.017	ng/u1	0.00
7) Phenol-d5	6.893	99	86	0.002	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	145	0.005	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	101	0.003	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	30	0.001	ng/u1	0.00
21) Nitrobenzene-d5	8.863	128	55	0.004	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	41	0.004	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	101	0.004	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.651	131	81	0.002	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	157	0.002	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	946	0.010	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	51	0.004	ng/u1	0.00
60) Fluorene-d10	15.363	176	1282	0.019	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	14	0.002	ng/u1	-0.02
73) Anthracene-d10	17.127	188	2037202	22.476	ng/u1	-0.11
81) Pyrene-d10	19.480	212	2169	0.023	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.445	264	2955	0.033	ng/u1	-0.01

Target Compounds	Qvalue

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

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