

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP011004.D
 Acq On : 18 Jul 2022 21:36
 Operator : CG/JU
 Sample : N3710-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 19 00:47:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 14 15:34:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	293986	20.000	ng/u1	0.00
20) Naphthalene-d8	10.475	136	1199736	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.340	164	648611	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.104	188	1313983	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.227	240	1288521	20.000	ng/u1	#-0.01
88) Perylene-d12	23.574	264	1324076	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	44599	5.469	ng/uL	0.00
4) Pyridine-d5	3.599	84	169045	8.020	ng/u1	0.00
7) Phenol-d5	6.858	99	192813	7.990	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	572170	36.872	ng/u1	0.00
11) 2-Chlorophenol-d4	7.216	132	575804	29.529	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	360138	18.641	ng/u1	-0.01
21) Nitrobenzene-d5	8.840	128	323173	36.963	ng/u1	0.00
24) 2-Nitrophenol-d4	9.563	143	303691	35.103	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	537233	33.956	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	577552	22.143	ng/u1	-0.01
46) Dimethylphthalate-d6	13.763	166	1939984	43.762	ng/u1	0.00
49) Acenaphthylene-d8	14.034	160	2143978	40.991	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	54550	6.080	ng/u1	0.00
60) Fluorene-d10	15.339	176	1618832	42.962	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	201847	31.044	ng/u1	-0.01
73) Anthracene-d10	17.204	188	2504470	44.834	ng/u1	-0.01
81) Pyrene-d10	19.463	212	2983309	44.084	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.421	264	2890896	45.458	ng/u1	-0.03

Target Compounds Qvalue

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

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