

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP011003.D
 Acq On : 18 Jul 2022 21:02
 Operator : CG/JU
 Sample : N3710-15
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0B41

Quant Time: Jul 19 00:47:16 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	322523	20.000	ng/u1	0.00
20) Naphthalene-d8	10.475	136	1350815	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.340	164	747390	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.104	188	1479175	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.227	240	1441766	20.000	ng/u1	#-0.01
88) Perylene-d12	23.580	264	1486561	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	43759	4.891	ng/uL	0.00
4) Pyridine-d5	3.599	84	161720	6.994	ng/u1	0.00
7) Phenol-d5	6.858	99	159595	6.028	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	500603	29.405	ng/u1	0.00
11) 2-Chlorophenol-d4	7.216	132	498276	23.292	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	306346	14.454	ng/u1	-0.01
21) Nitrobenzene-d5	8.840	128	286343	29.088	ng/u1	0.00
24) 2-Nitrophenol-d4	9.563	143	267171	27.428	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	460537	25.853	ng/u1	0.00
31) 4-Chloroaniline-d4	10.622	131	713616	24.300	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	1719221	33.656	ng/u1	0.00
49) Acenaphthylene-d8	14.034	160	1900886	31.540	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	46642	4.512	ng/u1	0.00
60) Fluorene-d10	15.340	176	1431210	32.962	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	186087	25.424	ng/u1	-0.01
73) Anthracene-d10	17.204	188	2328287	37.026	ng/u1	-0.01
81) Pyrene-d10	19.463	212	2794011	36.898	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.427	264	2711630	37.979	ng/u1	-0.02

Target Compounds	Qvalue

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

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