

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
 Data File : BP011071.D
 Acq On : 21 Jul 2022 20:13
 Operator : CG/JU
 Sample : N3709-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 21 23:25:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.669	152	370565	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.463	136	1505601	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.333	164	842217	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	1690602	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.221	240	1599155	20.000	ng/ul	# 0.00
88) Perylene-d12	23.574	264	1642859	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	43374	4.220	ng/uL	0.00
4) Pyridine-d5	3.593	84	191782	7.218	ng/ul	0.00
7) Phenol-d5	6.851	99	743987	24.458	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	490515	25.077	ng/ul	0.00
11) 2-Chlorophenol-d4	7.204	132	619174	25.191	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	522612	21.461	ng/ul	-0.01
21) Nitrobenzene-d5	8.834	128	293876	26.784	ng/ul	0.00
24) 2-Nitrophenol-d4	9.551	143	289310	26.648	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	487902	24.573	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.610	131	732078	22.366	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	1597853	27.759	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160	1935044	28.491	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143	180750	15.515	ng/ul	0.00
60) Fluorene-d10	15.333	176	1352338	27.639	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	121717	14.550	ng/ul	0.00
73) Anthracene-d10	17.198	188	2031891	28.271	ng/ul	0.00
81) Pyrene-d10	19.457	212	2404458	28.629	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.421	264	2274871	28.830	ng/ul	0.00

Target Compounds Qvalue

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

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