

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
 Data File : BP016238.D
 Acq On : 14 Jul 2023 21:56
 Operator : MA/JU
 Sample : 03457-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0C47

Quant Time: Jul 15 00:44:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 14 23:53:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	140644	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	682156	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	445726	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.380	188	1032154	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	1116710	20.000	ng/ul	0.01
88) Perylene-d12	23.968	264	1268469	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	10071	2.534	ng/uL	0.00
4) Pyridine-d5	3.752	84	93573	9.440	ng/ul	0.01
7) Phenol-d5	7.199	99	60150	4.735	ng/ul	0.05
9) Bis-(2-Chloroethyl)eth...	7.293	67	284604	33.131	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	235127	25.639	ng/ul	0.02
15) 4-Methylphenol-d8	8.722	113	155118	15.158	ng/ul	0.02
21) Nitrobenzene-d5	9.163	128	160647	34.135	ng/ul	0.01
24) 2-Nitrophenol-d4	9.892	143	174383	32.511	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.475	165	269804	29.047	ng/ul	0.03
31) 4-Chloroaniline-d4	10.981	131	368749	24.768	ng/ul	0.02
46) Dimethylphthalate-d6	14.033	166	1142958	33.616	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	1293745	34.721	ng/ul	0.00
54) 4-Nitrophenol-d4	14.933	143	9	0.001	ng/ul	0.00
60) Fluorene-d10	15.622	176	1007875	35.420	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.822	200	109579	18.391	ng/ul	0.03
73) Anthracene-d10	17.486	188	1672513	37.045	ng/ul	0.00
81) Pyrene-d10	19.715	212	2100933	36.257	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.798	264	2206079	36.215	ng/ul	0.00

Target Compounds Qvalue

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

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