

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
 Data File : BP017102.D
 Acq On : 11 Sep 2023 13:39
 Operator : MA/JU
 Sample : 04268-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0GH1

Quant Time: Sep 12 02:28:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 06:13:29 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	257949	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	1160462	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	725219	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.386	188	1578869	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	1346009	20.000	ng/ul	0.00
88) Perylene-d12	24.015	264	1441079	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	34216	5.063	ng/uL	0.00
4) Pyridine-d5	3.758	84	178300	8.749	ng/ul	0.00
7) Phenol-d5	7.110	99	127979	5.458	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	440220	27.993	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	358613	21.161	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	247714	13.355	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	231694	28.044	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	233999	27.442	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	406727	25.155	ng/ul	0.00
31) 4-Chloroaniline-d4	10.945	131	548615	21.160	ng/ul	0.00
46) Dimethylphthalate-d6	14.034	166	1651914	32.012	ng/ul	0.00
49) Acenaphthylene-d8	14.316	160	1679676	28.586	ng/ul	0.00
54) 4-Nitrophenol-d4	14.833	143	44872	4.215	ng/ul	0.00
60) Fluorene-d10	15.616	176	1388841	31.716	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	222275	24.946	ng/ul	0.00
73) Anthracene-d10	17.486	188	2303881	34.043	ng/ul	0.00
81) Pyrene-d10	19.727	212	2894093	40.829	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.851	264	2680597	38.748	ng/ul	0.00

Target Compounds Qvalue

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

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