

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018180.D
 Acq On : 15 Nov 2023 18:49
 Operator : MA/JU
 Sample : 05338-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDL2

Quant Time: Nov 15 22:00:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	65508	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	253506	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	149005	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	309749	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	275332	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	323453	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	6319	3.403	ng/uL	0.00
4) Pyridine-d5	3.611	84	21947	4.533	ng/ul	0.02
7) Phenol-d5	6.963	99	32326	5.111	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	93807	25.240	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	96798	19.328	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	63000	12.157	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	55445	24.276	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	61417	23.736	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	98702	21.402	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	90872	14.369	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	372401	27.087	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	379136	25.319	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	161	0.078	ng/ul	0.14
60) Fluorene-d10	15.469	176	300830	26.503	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	40465	18.071	ng/ul	0.00
73) Anthracene-d10	17.345	188	492326	29.388	ng/ul	0.00
81) Pyrene-d10	19.604	212	574301	31.227	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	585691	31.141	ng/ul	0.00

Target Compounds Qvalue

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

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