

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110624\  
 Data File : BM048481.D  
 Acq On : 06 Nov 2024 17:32  
 Operator : RC/JU  
 Sample : P4634-06  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

**Instrument :**  
 BNA\_M  
**ClientSampleId :**  
 CC0P1

**Manual Integrations**  
**APPROVED**  
 Reviewed By :Yogesh Patel 11/07/2024  
 Supervised By :mohammad ahmed 11/11/2024

Quant Time: Nov 06 22:22:10 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM110624.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 06 14:39:10 2024  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.678	152	3856	0.400	ng/ul	0.00
4) Naphthalene-d8	10.464	136	12162	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.326	164	6394	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.070	188	12581m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.274	240	9981m	0.400	ng/ul	0.01
23) Perylene-d12	23.487	264	8530m	0.400	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	18429	3.962	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.086	152	3806	0.242	ng/ul	0.01
18) Fluoranthene-d10	19.100	212	7694	0.275	ng/ul	0.00

Target Compounds Qvalue

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

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